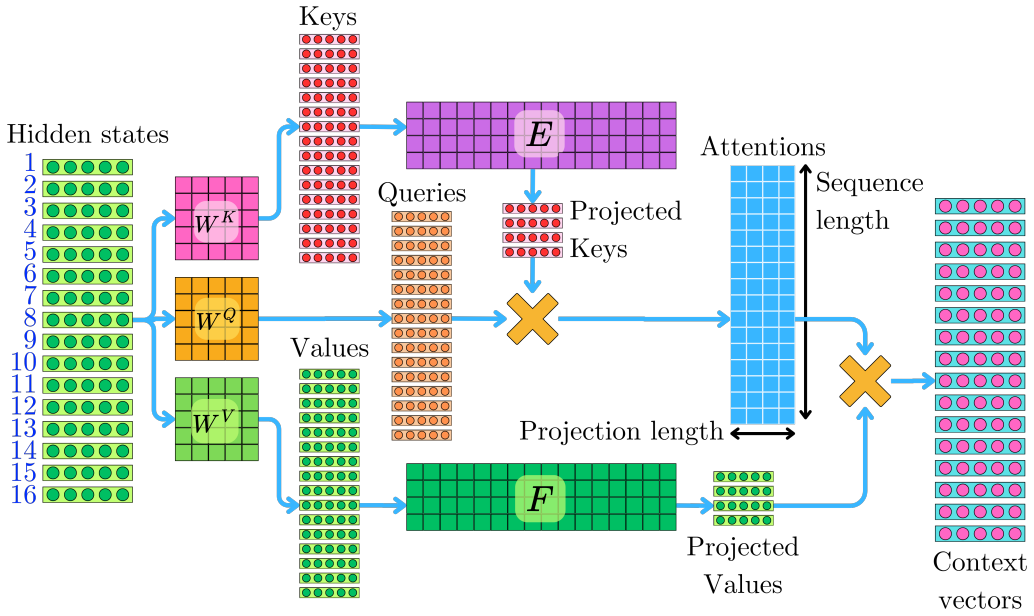


# Low-Rank Projection of Attention Matrices: Linformer

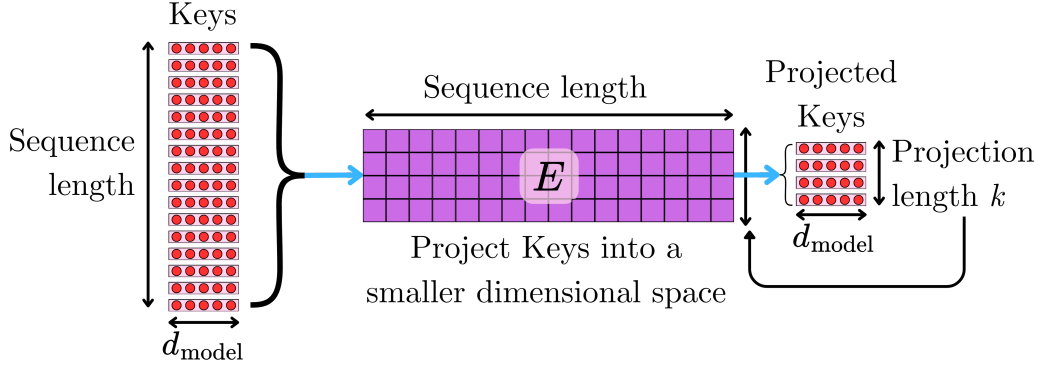
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Linear attention mechanisms represent a paradigm shift in transformer architecture by mathematically re-engineering the attention operation to achieve  $\mathcal{O}(n)$  complexity while maintaining global context awareness. Linformer introduced the idea that the token-token interaction matrix could be compressed into a smaller representation without too much information loss. Instead of computing the full  $N \times N$  interaction  $\frac{Q^\top K}{\sqrt{d_{\text{model}}}}$  (ignoring heads for simplicity), we could first project  $K$  into a lower rank dimension  $k$ , and compute the lower rank  $N \times k$  approximation:

$$\frac{Q^\top EK}{\sqrt{d_{\text{model}}}} \quad (1)$$

where  $E$  is a  $N \times k$  projection matrix that project  $K$  from the original dimension  $d_{\text{model}} \times N$  to  $d_{\text{model}} \times k$ . This leads to  $N \times k$  alignment score and attention matrices.



When we project with  $E$ , the approximation leads to the error:

$$\text{error} = \left| \frac{Q^\top K}{\sqrt{d_{\text{model}}}} - \frac{Q^\top EK}{\sqrt{d_{\text{model}}}} \right| \quad (2)$$

If the elements of  $E$  follow a Gaussian distribution  $\mathcal{N}(0, 1/k)$ , the Johnson–Lindenstrauss lemma guarantees that:

$$P[\text{error} > \epsilon] \leq e^{-\gamma \epsilon^2 k}. \quad (3)$$

This means that the probability that we choose  $E$  such that the error is greater than  $\epsilon$  is bounded by  $e^{-\gamma \epsilon^2 k}$ . If we choose  $k \rightarrow \infty$ , then  $P[\text{error} > \epsilon] \rightarrow 0$  for any  $\epsilon$ . A good choice is  $k \propto \log N / \epsilon^2$ , yielding:

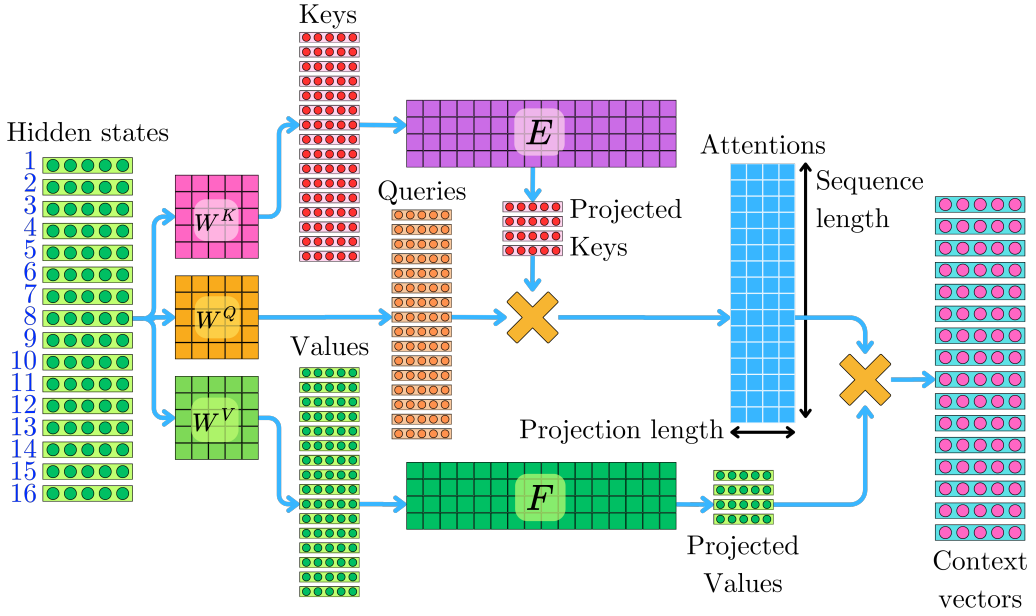
$$P[\text{error} > \epsilon] \leq N^{-\gamma}. \quad (4)$$

This means that we can choose an arbitrarily small  $\epsilon$  such that  $P[\text{error} > \epsilon] \rightarrow 0$  as the sequence length increases  $N \rightarrow \infty$ . Understand this as a mere theoretical guide that tells us that choosing  $k \propto \log N$  will guarantee smaller errors as  $N$  increases. In practice,  $k$  is chosen independently of  $N$ , leading to the  $\mathcal{O}(N)$  linear complexity while accepting the cost of the approximation error. Additionally,  $E$  is chosen as a parameter layer for the model to learn. For example, they showed that choosing  $k = 64$  with  $N = 512$  leads to slightly worse performance than the full attention.

Since the attention matrix has dimension  $N \times k$ , we also need to project the values:

$$C = \text{Softmax} \left( \frac{Q^\top EK}{\sqrt{d_{\text{model}}}} \right) FV \quad (5)$$

where  $F$  is the  $N \times k$  projection matrix for the tensor  $V$ . As for  $E$ ,  $F$  is also learned during training.



Projecting the keys and values  $EK$ ,  $FV$  leads to complexity  $\mathcal{O}(Nk)$ . Computing the alignment scores  $Q^\top EK$  and the context vectors  $C = AFV$  are also following  $\mathcal{O}(Nk)$ . Since we fix  $k$ , the overall time and space complexity is  $\mathcal{O}(N)$